## \*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

1358 REFERENCES IN FILE CA (1907 TO DATE)
63 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
1377 REFERENCES IN FILE CAPLUS (1907 TO DATE)

## => d his

(FILE 'HOME' ENTERED AT 10:24:35 ON 03 JAN 2008)

FILE 'REGISTRY' ENTERED AT 10:24:57 ON 03 JAN 2008 E 82657-04-3/RN

L1 1 S E3

FILE 'CAPLUS' ENTERED AT 10:25:23 ON 03 JAN 2008

L2 1377 S L1 ·

E WOOD+ALL/CT

L3 48 S L2 AND ("CONSTRUCTION MATERIALS" OR "WOOD" OR "LIGNOCELLULOSI

FILE 'REGISTRY' ENTERED AT 10:32:40 ON 03 JAN 2008

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L1
     ANSWER 1 OF 1 REGISTRY COPYRIGHT 2008 ACS on STN
RN
     82657-04-3 REGISTRY
ED
     Entered STN: 16 Nov 1984
     Cyclopropanecarboxylic acid, 3-[(1Z)-2-chloro-3,3,3-trifluoro-1-propen-1-
CN
     yl]-2,2-dimethyl-, (2-methyl[1,1'-biphenyl]-3-yl)methyl ester,
     (1R, 3R) -rel- (CA INDEX NAME)
OTHER CA INDEX NAMES:
     Cyclopropanecarboxylic acid, 3-(2-chloro-3,3,3-trifluoro-1-propenyl)-2,2-
     dimethyl-, (2-methyl[1,1'-biphenyl]-3-yl)methyl ester,
     [1\alpha, 3\alpha(Z)] - (\pm) -
CN
     Cyclopropanecarboxylic acid, 3-[(1Z)-2-chloro-3,3,3-trifluoro-1-propenyl]-
     2,2-dimethyl-, (2-methyl[1,1'-biphenyl]-3-yl)methyl ester, (1R,3R)-rel-
     (9CI)
OTHER NAMES:
     AGST 02002
CN
     Bifenthrin
CN
     Bifenthrine
CN
CN
     Biflex
CN
     Biflex FT
CN
     Biphenate
     Biphenthrin
CN
     Biphentrin
CN
     Brigade
CN
     Brigade 10WP
CN
CN
     Brigata Flo
CN
     Capture
CN
     Capture (pesticide)
CN
     Cyclopropanecarboxylic acid, 3-(2-chloro-3,3,3-trifluoro-1-propenyl)-2,2-
     dimethyl-, (2-methyl[1,1'-biphenyl]-3-yl)methyl ester,
     [1\alpha, 3\alpha(Z)] -
     Discipline
CN
CN
     Empower
     Fanfare
CN
     FMC 54800
CN
     Kiros EV
ÇN
CN
     Onyx
     Onyx (insecticide)
CN
CN
     Semafor
     Silencer
CN
     Talstar
CN
CN
     TalstarOne
FS
     STEREOSEARCH
     92880-79-0, 107497-60-9, 107538-32-9
DR
MF
     C23 H22 Cl F3 O2
CI
LC
                  AGRICOLA, ANABSTR, AQUIRE, BEILSTEIN*, BIOSIS, CA, CABA,
       CAPLUS, CASREACT, CBNB, CHEMCATS, CHEMLIST, CIN, CSCHEM, CSNB, EMBASE.
       HSDB*, IFICDB, IFIUDB, MEDLINE, MRCK*, PIRA, PROMT, RTECS*, TOXCENTER,
       USPAT2, USPATFULL
         (*File contains numerically searchable property data)
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Relative stereochemistry.

Double bond geometry as shown.